

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112316\
 Data File : BF091294.D
 Acq On : 24 Nov 2016 00:10
 Operator : UM/SJ
 Sample : H5740-10
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-04(0-0.16)

Manual Integrations
 APPROVED

sohil
 11/28/2016 5:18:11 PM

Quant Time: Nov 24 00:58:34 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 23 13:22:53 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	203683	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	806315	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	362733	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	489626	20.00	ng	0.00
75) Chrysene-d12	14.08	240	342188	20.00	ng	0.03
86) Perylene-d12	15.54	264	384458	20.00	ng	0.07

System Monitoring Compounds

5) 2-Fluorophenol	5.48	112	1164806	91.88	ng	0.01
7) Phenol-d6	6.51	99	1647293	101.14	ng	0.01
23) Nitrobenzene-d5	7.44	82	975928	69.68	ng	-0.01
41) 2,4,6-Tribromophenol	10.70	330	290701	79.47	ng	0.00
44) 2-Fluorobiphenyl	9.24	172	1268482	57.26	ng	0.00
78) Terphenyl-d14	13.00	244	906634	60.51	ng	0.01

Target Compounds

						Qvalue
10) Phenol	6.52	94	118153	6.50	ng	80
31) Naphthalene	8.19	128	177418	4.57	ng	98
32) Benzoic acid	7.87	122	40295	4.10	ng	95
37) 2-Methylnaphthalene	8.88	142	84836	3.19	ng	93
48) Acenaphthylene	9.78	152	1245394	40.57	ng	99
49) Dimethylphthalate	9.64	163	141260	6.24	ng	# 95
51) Acenaphthene	9.95	154	63033	3.02	ng	98
54) Dibenzofuran	10.12	168	122631	4.17	ng	# 92
57) Fluorene	10.46	166	106564m	4.74	ng	
68) Atrazine	11.10	200	17770	3.80	ng	85
70) Phenanthrene	11.42	178	935464	36.83	ng	98
71) Anthracene	11.49	178	2993144	115.25	ng	99
72) Carbazole	11.63	167	377510	16.33	ng	100
74) Fluoranthene	12.64	202	5363432	207.59	ng	93
77) Pyrene	12.88	202	6356273	251.40	ng	94
80) Benzo(a)anthracene	14.07	228	6941518	357.43	ng	# 90
82) Chrysene	14.11	228	4005596m	218.24	ng	
83) Bis(2-ethylhexyl)phthalate	14.04	149	67434	5.35	ng	# 87
85) Indeno(1,2,3-cd)pyrene	17.01	276	3203857	202.69	ng	# 92
87) Benzo(b)fluoranthene	15.14	252	8761091m	361.24	ng	
88) Benzo(k)fluoranthene	15.15	252	2652563m	133.91	ng	
89) Benzo(a)pyrene	15.51	252	4339515	210.32	ng	95
90) Dibenzo(a,h)anthracene	17.01	278	1170613	61.45	ng	# 87
91) Benzo(g,h,i)perylene	17.45	276	2652457	140.38	ng	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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