

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071917\
 Data File : BF096884.D
 Acq On : 19 Jul 2017 21:52
 Operator : SJ/JU
 Sample : I4280-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-071817-C

Manual Integrations
 APPROVED

Sohil
 7/20/2017 3:05:05 PM

Quant Time: Jul 20 05:16:47 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF071317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 18 13:33:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.59	152	104792	20.00	ng	-0.01
21) Naphthalene-d8	7.87	136	405714	20.00	ng	-0.02
38) Acenaphthene-d10	9.62	164	161072	20.00	ng	-0.02
63) Phenanthrene-d10	11.09	188	246215	20.00	ng	-0.01
75) Chrysene-d12	13.73	240	187447	20.00	ng	0.00
86) Perylene-d12	15.10	264	133240	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.19	112	554605	79.58	ng	0.00
7) Phenol-d6	6.24	99	670576	80.18	ng	-0.01
23) Nitrobenzene-d5	7.15	82	383758	58.59	ng	-0.02
41) 2,4,6-Tribromophenol	10.40	330	112875	82.10	ng	-0.02
44) 2-Fluorobiphenyl	8.95	172	671497	65.87	ng	-0.01
78) Terphenyl-d14	12.68	244	448445	41.49	ng	-0.01

Target Compounds

						Qvalue
10) Phenol	6.25	94	30431	3.219	ng	95
31) Naphthalene	7.89	128	176465	9.182	ng	100
37) 2-Methylnaphthalene	8.58	142	102838	8.393	ng	94
48) Acenaphthylene	9.47	152	94693	5.843	ng	97
49) Dimethylphthalate	9.35	163	95165	7.847	ng	98
51) Acenaphthene	9.65	154	51507	5.380	ng	98
54) Dibenzofuran	9.82	168	48989	3.651	ng	# 95
57) Fluorene	10.16	166	112242	11.321	ng	97
70) Phenanthrene	11.12	178	547448	40.892	ng	99
71) Anthracene	11.17	178	159928	11.815	ng	97
72) Carbazole	11.33	167	34448	2.628	ng	97
74) Fluoranthene	12.30	202	428049	33.404	ng	99
77) Pvrene	12.53	202	492897	28.974	ng	99
80) Benzo(a)anthracene	13.72	228	183287m	15.693	ng	
82) Chrysene	13.75	228	180852m	15.729	ng	
85) Indeno(1,2,3-cd)pyrene	16.39	276	50229	7.849	ng	96
87) Benzo(b)fluoranthene	14.73	252	130394m	16.226	ng	
88) Benzo(k)fluoranthene	14.74	252	45080m	5.924	ng	
89) Benzo(a)pvrene	15.04	252	110951	15.054	ng	# 93
90) Dibenzo(a,h)anthracene	16.39	278	15477	2.282	ng	# 62
91) Benzo(g,h,i)perylene	16.77	276	55608	7.657	ng	95

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

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