

Data Path : Z:\HPCHEM1\BNA F\DATA\BF081017\
 Data File : BF097583.D
 Acq On : 11 Aug 2017 2:58
 Operator : SJ/JU
 Sample : I4697-15 5X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NYSEG-PH4-ESMI-9B

Manual Integrations
 APPROVED

sohil
 8/14/2017 6:35:59 PM

Quant Time: Aug 11 07:46:27 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 07 16:02:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.39	152	110451	20.00	ng	-0.02
21) Naphthalene-d8	7.67	136	458240	20.00	ng	-0.02
38) Acenaphthene-d10	9.42	164	161791	20.00	ng	-0.02
63) Phenanthrene-d10	10.89	188	243649	20.00	ng	-0.02
75) Chrysene-d12	13.53	240	170944	20.00	ng	-0.01
86) Perylene-d12	14.86	264	169701	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.00	112	125492	17.13	ng	0.00
7) Phenol-d6	6.08	99	160520	19.19	ng	0.00
23) Nitrobenzene-d5	6.97	82	86148	11.03	ng	-0.02
41) 2,4,6-Tribromophenol	10.21	330	18479	14.46	ng	-0.02
44) 2-Fluorobiphenyl	8.75	172	120494	12.64	ng	-0.02
78) Terphenyl-d14	12.47	244	53081	6.33	ng	-0.02

Target Compounds

						Qvalue
31) Naphthalene	7.69	128	124764	5.776	ng	99
37) 2-Methylnaphthalene	8.39	142	50151	3.667	ng	97
48) Acenaphthylene	9.28	152	95146	6.096	ng	98
49) Dimethylphthalate	9.16	163	34113	2.777	ng	99
51) Acenaphthene	9.45	154	171592	18.663	ng	99
54) Dibenzofuran	9.62	168	165348	13.501	ng	98
57) Fluorene	9.96	166	268142	29.131	ng	96
70) Phenanthrene	10.92	178	1303360	99.838	ng	99
71) Anthracene	10.97	178	545334	40.785	ng	99
72) Carbazole	11.13	167	143191	10.889	ng	99
73) Di-n-butylphthalate	11.47	149	40125	2.468	ng	# 91
74) Fluoranthene	12.11	202	1198245	93.692	ng	99
77) Pvrene	12.33	202	1077586	77.000	ng	100
80) Benzo(a)anthracene	13.52	228	635202	57.428	ng	98
82) Chrysene	13.55	228	498832	46.186	ng	97
85) Indeno(1,2,3-cd)pyrene	16.05	276	178122	19.233	ng	96
87) Benzo(b)fluoranthene	14.52	252	582515m	57.103	ng	
88) Benzo(k)fluoranthene	14.54	252	196420m	20.206	ng	
89) Benzo(a)pvrene	14.82	252	435081	46.601	ng	96
90) Dibenzo(a,h)anthracene	16.04	278	46227m	6.038	ng	
91) Benzo(g,h,i)perylene	16.40	276	157294	19.591	ng	97

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

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