

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030518\
 Data File : BF103418.D
 Acq On : 5 Mar 2018 19:37
 Operator : SJ/JU
 Sample : J1722-19
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-17

Manual Integrations
 APPROVED

Sohil
 3/6/2018 4:23:13 PM

Quant Time: Mar 06 00:33:44 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 05 13:40:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	147763	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	613209	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	260715	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	378311	20.00	ng	0.00
75) Chrysene-d12	14.07	240	211890	20.00	ng	0.00
86) Perylene-d12	15.59	264	198617	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1020357	104.44	ng	0.00
7) Phenol-d6	6.52	99	1223704	102.36	ng	0.00
23) Nitrobenzene-d5	7.44	82	766709	71.40	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	177457	80.41	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1049199	65.60	ng	0.00
78) Terphenyl-d14	13.00	244	696386	79.00	ng	0.00
Target Compounds						
10) Phenol	6.53	94	93023	6.703	ng	Qvalue # 76
31) Naphthalene	8.17	128	64580	2.151	ng	98
48) Acenaphthylene	9.77	152	86104	3.238	ng	97
49) Dimethylphthalate	9.62	163	95806	4.581	ng	98
51) Acenaphthene	9.95	154	50387	3.249	ng	99
54) Dibenzofuran	10.12	168	63184	2.968	ng	97
57) Fluorene	10.46	166	74677	4.118	ng	95
70) Phenanthrene	11.44	178	1090872	52.396	ng	99
71) Anthracene	11.49	178	332058	15.554	ng	100
72) Carbazole	11.65	167	110331	5.190	ng	98
74) Fluoranthene	12.64	202	1488336	71.139	ng	98
77) Pvrene	12.87	202	1300899	78.746	ng	99
80) Benzo(a)anthracene	14.06	228	630198m	45.839	ng	
82) Chrysene	14.10	228	634273m	47.514	ng	
83) Bis(2-ethylhexyl)phthalate	14.02	149	25791	2.190	ng	# 94
85) Indeno(1,2,3-cd)pyrene	17.14	276	246318	23.307	ng	95
87) Benzo(b)fluoranthene	15.15	252	655302m	50.181	ng	
88) Benzo(k)fluoranthene	15.17	252	255552m	20.782	ng	
89) Benzo(a)pvrene	15.53	252	453988	38.930	ng	95
90) Dibenzo(a,h)anthracene	17.14	278	70005	7.769	ng	# 86
91) Benzo(g,h,i)perylene	17.62	276	229648	26.076	ng	96

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

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