

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF012318\
 Data File : BF102376.D
 Acq On : 24 Jan 2018 7:57
 Operator : SJ/JU
 Sample : J1214-01DL 4X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-PH3-MGP-01DL

Manual Integrations
 APPROVED

Sohil
 1/24/2018 2:47:27 PM

Quant Time: Jan 24 10:14:00 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 23 00:43:41 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	134229	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	557112	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	242958	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	368337	20.00	ng	0.00
75) Chrysene-d12	13.87	240	257315	20.00	ng	0.00
86) Perylene-d12	15.30	264	247325	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.32	112	209500	23.67	ng	0.00
7) Phenol-d6	6.35	99	257700	24.15	ng	-0.01
23) Nitrobenzene-d5	7.27	82	146587	15.12	ng	-0.01
41) 2,4,6-Tribromophenol	10.54	330	42334	17.59	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	191579	11.59	ng	-0.01
78) Terphenyl-d14	12.82	244	124526	10.91	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
31) Naphthalene	8.02	128	282221	10.146	ng	100
37) 2-Methylnaphthalene	8.71	142	53791	2.904	ng	99
48) Acenaphthylene	9.61	152	113027	4.296	ng	99
57) Fluorene	10.30	166	102730	6.235	ng	97
70) Phenanthrene	11.26	178	448086	20.876	ng	99
71) Anthracene	11.31	178	292530	13.353	ng	98
74) Fluoranthene	12.45	202	218366	9.977	ng	97
77) Pyrene	12.67	202	275748	13.860	ng	98
80) Benzo(a)anthracene	13.86	228	70515m	4.451	ng	
82) Chrysene	13.90	228	70468	4.380	ng	96
87) Benzo(b)fluoranthene	14.89	252	54846m	3.421	ng	
89) Benzo(a)pyrene	15.23	252	59350	4.105	ng	97
91) Benzo(g,h,i)perylene	17.09	276	25579	2.212	ng	92

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

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