

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040119\
 Data File : BF113498.D
 Acq On : 2 Apr 2019 9:21
 Operator : JU/SJ
 Sample : K2115-02DL 5X
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 QNWP8-MW30D-GWDL

Manual Integrations
 APPROVED

Sohil
 4/2/2019 3:50:04 PM

Quant Time: Apr 02 10:22:25 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	119396	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	324553	20.00	ng	0.02
39) Acenaphthene-d10	9.74	164	205139	20.00	ng	0.00
64) Phenanthrene-d10	11.22	188	350485	20.00	ng	0.00
76) Chrysene-d12	13.84	240	355016	20.00	ng	0.00
87) Perylene-d12	15.26	264	238531	20.00	ng	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.33	112	75271m	10.31	ng	0.01
7) Phenol-d6	6.36	99	59316	6.42	ng	0.00
23) Nitrobenzene-d5	7.27	82	120853	22.10	ng	0.00
42) 2,4,6-Tribromophenol	10.53	330	51302	20.25	ng	0.00
45) 2-Fluorobiphenyl	9.06	172	243520	1.22	ng	0.00
79) Terphenyl-d14	12.79	244	215306	13.37	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
6) Aniline	6.37	93	35390	3.146	ng	# 63
17) 2-Methylphenol	6.98	107	26335	3.963	ng	91
27) 2,4-Dimethylphenol	7.67	122	86433	19.922	ng	100
31) Naphthalene	8.04	128	11597169	731.216	ng	96
37) 2-Methylnaphthalene	8.71	142	1086100	109.749	ng	99
38) 1-Methylnaphthalene	8.80	142	1016190	106.757	ng	100
46) 1,1'-Biphenyl	9.16	154	264675	15.448	ng	98
52) Acenaphthene	9.77	154	462429	38.498	ng	99
55) Dibenzofuran	9.95	168	506559	28.052	ng	96
58) Fluorene	10.29	166	198447	14.146	ng	97
71) Phenanthrene	11.24	178	213093	12.243	ng	96
73) Carbazole	11.45	167	209076	12.394	ng	98

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

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