

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE050119\
 Data File : BE099461.D
 Acq On : 1 May 2019 13:16
 Operator : JU/SJ
 Sample : K2589-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PST-001-B197-042519

Quant Time: May 02 08:10:19 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE042919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 29 18:36:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	1810	0.40	ng	-0.01
7) Naphthalene-d8	10.60	136	6600	0.40	ng	0.00
13) Acenaphthene-d10	14.46	164	5305	0.40	ng	0.00
19) Phenanthrene-d10	17.19	188	18865	0.40	ng	0.00
27) Chrysene-d12	21.37	240	24016	0.40	ng	0.00
34) Perylene-d12	23.86	264	27666	0.40	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.39	112	2137	0.43	ng	0.00
5) Phenol-d6	6.97	99	2867	0.43	ng	0.00
8) Nitrobenzene-d5	8.96	82	2359	0.38	ng	0.00
11) 2-Methylnaphthalene-d10	12.19	152	4487	0.38	ng	0.00
14) 2,4,6-Tribromophenol	15.93	330	1530	0.46	ng	-0.01
15) 2-Fluorobiphenyl	13.07	172	6766	0.33	ng	-0.02
25) Fluoranthene-d10	19.22	212	85492	0.34	ng	0.00
29) Terphenyl-d14	19.81	244	16821	0.38	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
23) Phenanthrene	17.23	178	2080	0.043	ng	# 94
26) Fluoranthene	19.25	202	6583	0.090	ng	# 99
28) Pyrene	19.61	202	6328	0.101	ng	# 96
30) Benzo(a)anthracene	21.34	228	3774	0.049	ng	# 88
31) Chrysene	21.40	228	3904	0.052	ng	# 92
32) Bis(2-ethylhexyl)phthalate	21.26	149	15388	0.071	ng	# 98
33) Indeno(1,2,3-cd)pyrene	26.50	276	2125	0.024	ng	# 91
35) Benzo(b)fluoranthene	23.09	252	4893	0.061	ng	# 48
37) Benzo(a)pyrene	23.75	252	3165	0.042	ng	# 52
39) Benzo(g,h,i)perylene	27.32	276	2123	0.027	ng	# 68

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

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