

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE112219\
 Data File : BE100759.D
 Acq On : 22 Nov 2019 15:10
 Operator : JU
 Sample : K5969-05
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 HF-102S

Quant Time: Nov 22 16:52:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE112119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 22 06:40:44 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	5894	0.40	ng	0.00
7) Naphthalene-d8	10.55	136	25449	0.40	ng	0.00
13) Acenaphthene-d10	14.40	164	13983	0.40	ng	0.00
19) Phenanthrene-d10	17.13	188	32645	0.40	ng	0.00
27) Chrysene-d12	21.32	240	35001	0.40	ng	0.00
34) Perylene-d12	23.77	264	39166	0.40	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.37	112	2397	0.15	ng	0.00
5) Phenol-d6	6.93	99	3049	0.13	ng	0.00
8) Nitrobenzene-d5	8.90	82	4253	0.52	ng	0.00
11) 2-Methylnaphthalene-d10	12.14	152	13991	0.37	ng	0.00
14) 2,4,6-Tribromophenol	15.88	330	1061	0.54	ng	0.00
15) 2-Fluorobiphenyl	13.03	172	21375	0.39	ng	0.00
25) Fluoranthene-d10	19.17	212	163121	0.41	ng	0.00
29) Terphenyl-d14	19.77	244	38725	0.48	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.35	88	32884	2.974	ng	99
26) Fluoranthene	19.20	202	3062	0.025	ng	# 89
28) Pyrene	19.56	202	2933	0.027	ng	# 88
30) Benzo(a)anthracene	21.29	228	2683	0.023	ng	# 80
31) Chrysene	21.35	228	2443	0.020	ng	# 85
32) Bis(2-ethylhexyl)phthalate	21.26	149	12254	0.023	ng	# 98
33) Indeno(1,2,3-cd)pyrene	26.35	276	1963	0.053	ng	97
39) Benzo(g,h,i)perylene	27.14	276	1801	0.037	ng	# 74

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

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