

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE032319\
 Data File : BE099104.D
 Acq On : 23 Mar 2019 10:08
 Operator : JU/SJ
 Sample : K1947-09
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PST-025-B033-031319

Manual Integrations
 APPROVED

Sohil
 3/25/2019 12:36:53 PM

Quant Time: Mar 24 02:56:21 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE032219.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Sat Mar 23 04:40:11 2019

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	3578	0.40	ng	0.00
7) Naphthalene-d8	10.63	136	14108	0.40	ng	-0.01
13) Acenaphthene-d10	14.49	164	8806	0.40	ng	0.00
19) Phenanthrene-d10	17.22	188	23363	0.40	ng	0.00
27) Chrysene-d12	21.39	240	35566	0.40	ng	0.00
34) Perylene-d12	23.89	264	39503	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.41	112	4966	0.47	ng	0.00
5) Phenol-d6	6.99	99	6840	0.45	ng	0.00
8) Nitrobenzene-d5	8.99	82	4111	0.67	ng	0.00
11) 2-Methylnaphthalene-d10	12.22	152	9451	0.37	ng	-0.01
14) 2,4,6-Tribromophenol	15.96	330	1381	0.53	ng	0.00
15) 2-Fluorobiphenyl	13.10	172	14381	0.41	ng	-0.01
25) Fluoranthene-d10	19.24	212	117168	0.38	ng	0.00
29) Terphenyl-d14	19.84	244	23228	0.31	ng	0.00
Target Compounds						Qvalue
22) Pentachlorophenol	16.85	266	422	0.337	ng	# 82
23) Phenanthrene	17.26	178	7565	0.109	ng	98
26) Fluoranthene	19.27	202	14434	0.148	ng	97
28) Pyrene	19.63	202	14454	0.120	ng	95
30) Benzo(a)anthracene	21.37	228	9291	0.074	ng	99
31) Chrysene	21.43	228	9153	0.073	ng	96
32) Bis(2-ethylhexyl)phthalate	21.29	149	22783	0.140	ng	98
33) Indeno(1,2,3-cd)pyrene	26.55	276	4667	0.035	ng	# 88
35) Benzo(b)fluoranthene	23.13	252	11011	0.081	ng	# 80
36) Benzo(k)fluoranthene	23.17	252	4897m	0.037	ng	
37) Benzo(a)pyrene	23.78	252	7878	0.062	ng	# 89
39) Benzo(a,h,i)perylene	27.37	276	4866	0.039	ng	# 93

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

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