

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE032319\
 Data File : BE099116.D
 Acq On : 23 Mar 2019 17:39
 Operator : JU/SJ
 Sample : K2014-12
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PST-026-B121-031919

Manual Integrations
 APPROVED

Sohil
 3/25/2019 12:37:05 PM

Quant Time: Mar 24 03:28:13 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	3538	0.40	ng	0.00
7) Naphthalene-d8	10.63	136	13950	0.40	ng	-0.01
13) Acenaphthene-d10	14.49	164	9216	0.40	ng	0.00
19) Phenanthrene-d10	17.21	188	26396	0.40	ng	0.00
27) Chrysene-d12	21.38	240	38595	0.40	ng	-0.01
34) Perylene-d12	23.90	264	41998	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.41	112	4222	0.40	ng	0.00
5) Phenol-d6	6.99	99	6175	0.41	ng	0.00
8) Nitrobenzene-d5	8.99	82	4061	0.67	ng	-0.01
11) 2-Methylnaphthalene-d10	12.22	152	8810	0.35	ng	-0.01
14) 2,4,6-Tribromophenol	15.96	330	1583	0.58	ng	0.00
15) 2-Fluorobiphenyl	13.10	172	13464	0.37	ng	-0.01
25) Fluoranthene-d10	19.24	212	133979	0.38	ng	0.00
29) Terphenyl-d14	19.83	244	26235	0.32	ng	0.00
Target Compounds						
16) Acenaphthylene	14.20	152	1101	0.022	ng	Qvalue 96
23) Phenanthrene	17.25	178	8839	0.113	ng	99
26) Fluoranthene	19.27	202	16406	0.149	ng	# 96
28) Pyrene	19.63	202	20230	0.155	ng	96
30) Benzo(a)anthracene	21.37	228	10197	0.075	ng	94
31) Chrysene	21.42	228	11023	0.081	ng	94
32) Bis(2-ethylhexyl)phthalate	21.29	149	24245	0.137	ng	99
33) Indeno(1,2,3-cd)pyrene	26.54	276	6621	0.045	ng	97
35) Benzo(b)fluoranthene	23.12	252	13493	0.093	ng	# 84
36) Benzo(k)fluoranthene	23.17	252	4478m	0.031	ng	
37) Benzo(a)pyrene	23.78	252	9814	0.073	ng	# 89
39) Benzo(a,h,i)perylene	27.37	276	7451	0.056	ng	96

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

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