

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042618\
 Data File : BE096177.D
 Acq On : 26 Apr 2018 12:46
 Operator : SJ/JU
 Sample : J2431-11DL 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 DASP2DL

Manual Integrations
 APPROVED

Sohil
 4/27/2018 7:28:44 PM

Quant Time: Apr 27 02:53:52 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE042418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 27 02:29:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	672	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	2683	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	1649	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	4151	0.40	ng/ul	0.00
16) Chrysene-d12	21.75	240	4654	0.40	ng/ul	0.00
20) Perylene-d12	24.35	264	4119	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	261	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	874	0.05	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.23	128	13085	1.752	ng/ul#	88
5) 2-Methylnaphthalene	12.81	142	5565	1.112	ng/ul	100
7) Acenaphthylene	14.68	152	2884	0.339	ng/ul	98
8) Acenaphthene	15.01	153	11563	1.887	ng/ul	94
9) Fluorene	15.97	166	15435	1.903	ng/ul#	90
11) Pentachlorophenol	17.31	266	643	0.644	ng/ul	96
12) Phenanthrene	17.70	178	87000	7.119	ng/ul#	88
13) Anthracene	17.79	178	20095	1.639	ng/ul#	92
15) Fluoranthene	19.67	202	111200	5.953	ng/ul	82
17) Pyrene	20.02	202	48036	5.145	ng/ul#	77
18) Benzo(a)anthracene	21.73	228	26849	1.906	ng/ul#	86
19) Chrysene	21.79	228	41871	3.234	ng/ul	97
21) Benzo(b)fluoranthene	23.55	252	92155m	5.909	ng/ul	
22) Benzo(k)fluoranthene	23.59	252	20015m	1.342	ng/ul	
23) Benzo(a)pyrene	24.24	252	53893	3.977	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	27.13	276	29257	1.603	ng/ul#	73
25) Dibenzo(a,h)anthracene	27.13	278	7496	0.490	ng/ul	91
26) Benzo(g,h,i)perylene	28.00	276	21216	1.335	ng/ul#	91

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

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