

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042618\
 Data File : BE096176.D
 Acq On : 26 Apr 2018 12:10
 Operator : SJ/JU
 Sample : J2431-13DL 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 DASP4DL

Manual Integrations
 APPROVED

Sohil
 4/27/2018 7:28:42 PM

Quant Time: Apr 27 02:49:58 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	732	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	2837	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	1657	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	4177	0.40	ng/ul	0.00
16) Chrysene-d12	21.75	240	5359	0.40	ng/ul	0.00
20) Perylene-d12	24.35	264	3927	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	214	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	612	0.03	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.23	128	3021	0.382	ng/ul#	89
5) 2-Methylnaphthalene	12.81	142	5023	0.949	ng/ul	99
7) Acenaphthylene	14.68	152	3487	0.408	ng/ul#	94
8) Acenaphthene	15.01	153	28398	4.611	ng/ul	93
9) Fluorene	15.98	166	83573	10.256	ng/ul#	91
12) Phenanthrene	17.70	178	990860	80.571	ng/ul#	87
13) Anthracene	17.79	178	208989	16.939	ng/ul#	90
15) Fluoranthene	19.68	202	1121847	59.679	ng/ul	81
17) Pyrene	20.03	202	584130	54.333	ng/ul#	76
18) Benzo(a)anthracene	21.73	228	260218	16.039	ng/ul#	84
19) Chrysene	21.79	228	181407	12.168	ng/ul	97
21) Benzo(b)fluoranthene	23.55	252	140231m	9.431	ng/ul	
22) Benzo(k)fluoranthene	23.60	252	47725m	3.357	ng/ul	
23) Benzo(a)pyrene	24.24	252	85445	6.614	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	27.14	276	30834	1.771	ng/ul#	74
25) Dibenzo(a,h)anthracene	27.14	278	8863	0.608	ng/ul	91
26) Benzo(a,h,i)perylene	28.00	276	22306	1.472	ng/ul#	92

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

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