

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020518\
 Data File : BE095516.D
 Acq On : 6 Feb 2018 6:19
 Operator : SJ/JU
 Sample : J1315-04
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C61

Manual Integrations
 APPROVED

Sohil
 2/6/2018 2:31:52 PM

Quant Time: Feb 06 07:57:14 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 06 04:22:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1893	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	5458	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	3541	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8153	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	8276	0.40	ng/ul	0.00
20) Perylene-d12	24.43	264	7514m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	2824	0.32	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	8566	0.27	ng/ul	0.00
Target Compounds				Qvalue		
11) Pentachlorophenol	17.36	266	1352	0.856	ng/ul	97
13) Anthracene	17.85	178	745m	0.030	ng/ul	
15) Fluoranthene	19.72	202	2711	0.076	ng/ul	85
17) Pyrene	20.07	202	2728m	0.098	ng/ul	
18) Benzo(a)anthracene	21.78	228	973	0.038	ng/ul#	82
19) Chrysene	21.84	228	1140	0.042	ng/ul#	79
21) Benzo(b)fluoranthene	23.62	252	2098m	0.079	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.24	276	584m	0.022	ng/ul	
26) Benzo(g,h,i)perylene	28.11	276	553m	0.025	ng/ul	

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

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