

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091919\
 Data File : BP000324.D
 Acq On : 19 Sep 2019 12:53
 Operator : HP/JU
 Sample : K4878-22
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TPA-9

Quant Time: Sep 20 04:42:34 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP091619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 19 14:44:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	215755	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	791726	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	449722	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	852013	20.00	ng	0.00
76) Chrysene-d12	13.99	240	757346	20.00	ng	0.00
87) Perylene-d12	15.47	264	831302	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1460717	116.88	ng	0.02
7) Phenol-d6	6.43	99	1695785	110.69	ng	0.00
23) Nitrobenzene-d5	7.35	82	1176829	95.90	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	766940	171.94	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	2600606	104.08	ng	0.00
79) Terphenyl-d14	12.93	244	3285408	91.01	ng	0.00
Target Compounds						
71) Phenanthrene	11.35	178	90891	2.129	ng	98
73) Carbazole	11.56	167	87111	2.201	ng	96

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

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