

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091819\
 Data File : BP000295.D
 Acq On : 18 Sep 2019 14:24
 Operator : HP/JU
 Sample : K4925-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SS-3

Quant Time: Sep 19 03:44:08 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP091619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 16 18:59:05 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	296881	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	1080315	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	605680	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	1123656	20.00	ng	0.00
76) Chrysene-d12	13.99	240	930957	20.00	ng	0.00
87) Perylene-d12	15.47	264	984119	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1517220	88.23	ng	0.01
7) Phenol-d6	6.43	99	2020965	95.87	ng	0.00
23) Nitrobenzene-d5	7.35	82	1160043	69.28	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	586850	97.69	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	2336753	69.44	ng	0.00
79) Terphenyl-d14	12.93	244	2633596	59.35	ng	0.00
Target Compounds						
10) Phenol	6.43	94	77298	3.227	ng	85
50) Dimethylphthalate	9.55	163	323657	7.843	ng	98
75) Fluoranthene	12.55	202	204851	3.388	ng	95
78) Pyrene	12.78	202	179361	2.575	ng	99

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

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