

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP091619\
 Data File : BP000251.D
 Acq On : 16 Sep 2019 21:12
 Operator : HP/JU
 Sample : K4871-08
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 VE-3-1

Quant Time: Sep 17 12:08:13 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	339378	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	1157962	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	588672	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	988405	20.00	ng	0.00
76) Chrysene-d12	13.99	240	915753	20.00	ng	0.00
87) Perylene-d12	15.47	264	1033858	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	836266	42.54	ng	0.00
7) Phenol-d6	6.41	99	661690	27.46	ng	-0.01
23) Nitrobenzene-d5	7.35	82	1377534	76.75	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	628165	107.58	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	2633171	80.51	ng	0.00
79) Terphenyl-d14	12.93	244	2979528	68.26	ng	0.00
Target Compounds						
10) Phenol	6.42	94	153691	5.614	ng	100
50) Dimethylphthalate	9.55	163	250080	6.235	ng	# 80
52) Acenaphthene	9.87	154	161609	4.987	ng	94
58) Fluorene	10.39	166	158801	4.488	ng	# 88
71) Phenanthrene	11.36	178	355032	7.168	ng	# 89

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

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