

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP092719\
 Data File : BP000455.D
 Acq On : 27 Sep 2019 14:19
 Operator : JU
 Sample : K4663-09
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 OK-02-092419

Quant Time: Sep 27 23:35:22 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	233718	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	848422	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	468689	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	876870	20.00	ng	0.00
76) Chrysene-d12	14.00	240	787602	20.00	ng	0.00
87) Perylene-d12	15.49	264	858047	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1760738	130.06	ng	0.02
7) Phenol-d6	6.43	99	2200869	132.62	ng	0.01
23) Nitrobenzene-d5	7.35	82	1272121	96.74	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	733791	157.85	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	2681508	102.97	ng	0.00
79) Terphenyl-d14	12.94	244	3436909	91.54	ng	0.00
Target Compounds						
10) Phenol	6.45	94	132661	7.036	ng	97
50) Dimethylphthalate	9.55	163	380689	11.922	ng	99

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

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