

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081519\
 Data File : BF116286.D
 Acq On : 15 Aug 2019 18:56
 Operator : HP/JU
 Sample : K4333-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FRAC-TANK

Manual Integrations
 APPROVED

Jagrut
 8/16/2019 5:34:08 PM

Quant Time: Aug 16 11:12:36 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 11:13:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	115410	20.00	ng	-0.03
21) Naphthalene-d8	8.10	136	463882	20.00	ng	-0.03
39) Acenaphthene-d10	9.85	164	254243	20.00	ng	-0.04
64) Phenanthrene-d10	11.34	188	444643	20.00	ng	-0.03
76) Chrysene-d12	13.99	240	286294	20.00	ng	-0.02
87) Perylene-d12	15.44	264	313666	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	391151	56.74	ng	0.03
7) Phenol-d6	6.49	99	320348	36.83	ng	0.00
23) Nitrobenzene-d5	7.39	82	695246	86.51	ng	-0.03
42) 2,4,6-Tribromophenol	10.66	330	323659	131.30	ng	-0.02
45) 2-Fluorobiphenyl	9.17	172	1203981	88.70	ng	-0.03
79) Terphenyl-d14	12.93	244	1313510	96.68	ng	-0.02
Target Compounds						
10) Phenol	6.50	94	107274	10.473	ng	92
20) 3+4-Methylphenols	7.25	107	180909	23.683	ng	# 60
27) 2,4-Dimethylphenol	7.79	122	12537	2.037	ng	87
32) Benzoic acid	8.04	122	492065m	113.414	ng	
43) 2,4,6-Trichlorophenol	9.10	196	12909	2.759	ng	98
50) Dimethylphthalate	9.57	163	158630	9.163	ng	# 79
70) Pentachlorophenol	11.17	266	7477	2.798	ng	93
73) Carbazole	11.59	167	77602	3.718	ng	98

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

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