

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080721\
 Data File : BP006571.D
 Acq On : 07 Aug 2021 16:40
 Operator : CG/JU
 Sample : M3163-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGF01

Quant Time: Aug 08 02:35:44 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Aug 07 13:59:25 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.752	152	166656	20.000	ng/u1	0.00
20) Naphthalene-d8	10.534	136	716944	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.381	164	395520	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.127	188	785670	20.000	ng/u1	0.00
79) Chrysene-d12	21.227	240	775144	20.000	ng/u1	0.00
88) Perylene-d12	23.551	264	794923	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	43572	9.759	ng/uL	0.00
4) Pyridine-d5	3.681	84	617542	46.592	ng/u1	0.00
7) Phenol-d5	6.928	99	715863	45.736	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.099	67	477626	49.043	ng/u1	0.00
11) 2-Chlorophenol-d4	7.287	132	556047	47.817	ng/u1	0.00
15) 4-Methylphenol-d8	8.457	113	566698	46.040	ng/u1	0.00
21) Nitrobenzene-d5	8.904	128	273608	51.253	ng/u1	0.00
24) 2-Nitrophenol-d4	9.622	143	278519	54.642	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.157	165	442946	44.346	ng/u1	0.00
31) 4-Chloroaniline-d4	10.675	131	723190	43.979	ng/u1	0.00
46) Dimethylphthalate-d6	13.798	166	1411115	50.128	ng/u1	0.00
49) Acenaphthylene-d8	14.075	160	1764519	51.020	ng/u1	0.00
54) 4-Nitrophenol-d4	14.586	143	228863	40.946	ng/u1	0.00
60) Fluorene-d10	15.375	176	1167448	49.333	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.498	200	160518	36.648	ng/u1	0.00
73) Anthracene-d10	17.233	188	1758991	49.596	ng/u1	0.00
81) Pyrene-d10	19.474	212	1948515	48.872	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.404	264	1949782	48.267	ng/u1	0.00

Target Compounds	Qvalue

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

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