

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072821\
 Data File : BP006402.D
 Acq On : 28 Jul 2021 22:04
 Operator : CG/JU
 Sample : M3107-09
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGCZ9

Quant Time: Jul 29 01:08:25 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 24 04:29:12 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	356539	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.563	136	1564488	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.410	164	919540	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.157	188	1754171	20.000	ng/u1	-0.01
79) Chrysene-d12	21.257	240	1484958	20.000	ng/u1	-0.01
88) Perylene-d12	23.598	264	1304129	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	67831	7.102	ng/uL	0.00
4) Pyridine-d5	3.699	84	975618	34.407	ng/u1	0.00
7) Phenol-d5	6.958	99	1216898	36.341	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	800137	38.403	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	941973	37.864	ng/u1	-0.01
15) 4-Methylphenol-d8	8.487	113	989689	37.583	ng/u1	-0.01
21) Nitrobenzene-d5	8.940	128	468844	40.247	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.652	143	483670	43.484	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.187	165	808425	37.090	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.704	131	1269835	35.387	ng/u1	-0.01
46) Dimethylphthalate-d6	13.828	166	2578099	39.393	ng/u1	-0.01
49) Acenaphthylene-d8	14.098	160	3170779	39.435	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.610	143	417238	32.108	ng/u1	0.00
60) Fluorene-d10	15.404	176	2085085	37.898	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.522	200	292316	29.892	ng/u1	-0.01
73) Anthracene-d10	17.257	188	3118364	39.380	ng/u1	-0.01
81) Pyrene-d10	19.498	212	3301178	43.221	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.445	264	2584559	38.999	ng/u1	-0.02

Target Compounds	Qvalue

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

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