

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072821\
 Data File : BP006393.D
 Acq On : 28 Jul 2021 16:56
 Operator : CG/JU
 Sample : PB137929BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK929

Quant Time: Jul 28 17:24:35 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 28 15:50:13 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	356452	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.569	136	1573991	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.410	164	926377	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.163	188	1814953	20.000	ng/u1	0.00
79) Chrysene-d12	21.263	240	1663781	20.000	ng/u1	0.00
88) Perylene-d12	23.610	264	1498565	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	63954	6.697	ng/uL	0.00
4) Pyridine-d5	3.699	84	925669	32.653	ng/u1	0.00
7) Phenol-d5	6.958	99	1085036	32.411	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	705651	33.877	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	837206	33.661	ng/u1	-0.01
15) 4-Methylphenol-d8	8.487	113	870498	33.065	ng/u1	-0.01
21) Nitrobenzene-d5	8.940	128	417951	35.662	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.657	143	429909	38.417	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.187	165	703332	32.074	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.710	131	1171708	32.456	ng/u1	0.00
46) Dimethylphthalate-d6	13.834	166	2322038	35.219	ng/u1	0.00
49) Acenaphthylene-d8	14.104	160	2840531	35.067	ng/u1	0.00
54) 4-Nitrophenol-d4	14.610	143	412131	31.481	ng/u1	0.00
60) Fluorene-d10	15.404	176	1887836	34.060	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.522	200	282040	27.875	ng/u1	-0.01
73) Anthracene-d10	17.263	188	2997838	36.590	ng/u1	0.00
81) Pyrene-d10	19.504	212	3222570	37.657	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.457	264	2766611	36.330	ng/u1	0.00

Target Compounds Qvalue

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

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