

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
 Data File : BP007618.D
 Acq On : 23 Oct 2021 08:12
 Operator : CG/JU
 Sample : PB140107BL
 Misc :
 ALS Vial : 78 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK107

Quant Time: Oct 25 01:00:03 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 20 10:49:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	234612	20.000	ng/u1	0.00
20) Naphthalene-d8	10.746	136	954205	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.581	164	629903	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.334	188	1344572	20.000	ng/u1	0.00
79) Chrysene-d12	21.422	240	1414575	20.000	ng/u1	0.00
88) Perylene-d12	23.869	264	1451663	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	34575	5.821	ng/uL	0.00
4) Pyridine-d5	3.787	84	529725	32.577	ng/u1	0.00
7) Phenol-d5	7.111	99	681544	37.198	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	364881	33.390	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.475	132	541084	37.540	ng/u1	0.00
15) 4-Methylphenol-d8	8.652	113	537494	38.085	ng/u1	0.00
21) Nitrobenzene-d5	9.099	128	254511	42.147	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.828	143	286080	50.719	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	547369	39.786	ng/u1	0.00
31) 4-Chloroaniline-d4	10.881	131	713302	38.341	ng/u1	0.00
46) Dimethylphthalate-d6	13.993	166	1634205	37.500	ng/u1	0.00
49) Acenaphthylene-d8	14.275	160	1967498	35.925	ng/u1	0.00
54) 4-Nitrophenol-d4	14.775	143	276252	39.506	ng/u1	0.00
60) Fluorene-d10	15.575	176	1388345	38.069	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	93378	17.332	ng/u1	0.00
73) Anthracene-d10	17.434	188	2277296	39.881	ng/u1	0.00
81) Pyrene-d10	19.669	212	2690639	38.048	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.710	264	2770866	38.361	ng/u1	0.00

Target Compounds Qvalue

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

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